## **CLAIMS**

1. A compound of the following formula or a pharmaceutically acceptable salt thereof:

wherein:

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A represents O, NH, or CH<sub>2</sub>;

 $R_1$  represents a single bond, a lower alkyl group, a lower alkenyl group, a lower alkynyl group, or a formula of  $Y_1$ -W, wherein  $Y_1$  represents a lower alkyl group, a lower alkenyl group or dioxanyl; W represents a single bond or an oxygen atom; said lower alkyl group, lower alkenyl group, or lower alkynyl group may be substituted with one or more substituent(s), which is/are the same or different, selected from the group consisting of <substituent group  $\beta$ >;

 $R_2$  represents a phenyl group, a naphthyl group, or a five- or six-membered aromatic or aliphatic heterocyclic ring having at least one atom selected from N, S, or O, selected from the group consisting of <substituent group  $\alpha$ >, wherein said phenyl group, naphthyl group, aromatic or aliphatic heterocyclic ring may be substituted with one or more substituent(s), which is/are the same or different, selected from the group consisting of <substituent group  $\beta$ > and/or a lower alkyl group substituted with one or more substituent(s), which is/are the same or different, selected from the group of <substituent group  $\beta$ >; when A represents an oxygen atom,  $R_2$  may represent hydrogen atom; with a proviso that A is NH and  $R_1$  is CH<sub>2</sub>,  $R_2$  is not any of a substituted phenyl group, a hydroxymethyl-substituted naphthyl group, an unsubstituted thienyl group, a hydroxymethyl-substituted pyridyl group, an unsubstituted thienyl group, a hydroxymethyl-substituted 2-thienyl group, a

mono(hydroxymethyl)-substituted 3-thienyl group, an unsubstituted furyl group and a hydroxymethyl-substituted furyl group;

G represents a hexose group or a pentose group; <substituent group  $\alpha$ > represents the following:

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and;  $\langle$ substituent group  $\beta \rangle$  represents the following:

a hydroxyl group, a cyano group, a halogen atom, a nitro group, a carboxyl group, a carbamoyl group, a formyl group, a lower alkanoyl group, a lower alkanoyl group, a lower alkoxy group, a hydroxyl lower alkoxy group, a lower alkoxy carbonyl group, a lower alkyl carbamoyl group, a di-lower alkyl carbamoyloxy group, a di-lower alkyl carbamoyloxy group, a namino group, a lower alkyl amino group, a di-lower alkyl amino group, a tri-lower alkyl ammonio group, a lower alkanoyl

amino group, an aroyl amino group, a lower alkanoyl amidino group, a hydroxyl imino group, a lower alkoxy imino group, a lower alkyl thio group, a lower alkyl sulfinyl group, a lower alkyl sulfonyl group, a lower alkyl sulfonylamino group and a sulfamoyl group.

- 5 2. The compound according to claim 1 or a pharmaceutically acceptable salt thereof, wherein G represents β-glucopyranosyl group; the positions of substitution of the hydroxyl groups on the indolopyrrolocarbazole ring are the 2- and 10-positions; R<sub>1</sub> represents a lower alkyl group; and R<sub>2</sub> represents a five- or six-membered aromatic or aliphatic heterocyclic ring having at least one atom selected from N, S, or O, selected from the group consisting of the <substituent group α>.
  - 3. The compound according to claim 2 or a pharmaceutically acceptable salt thereof, wherein the <substituent group  $\alpha>$  represents the following:

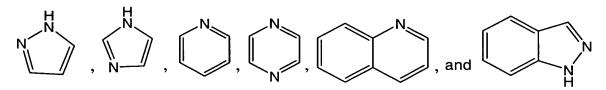
$$\begin{array}{c} \stackrel{H}{\longrightarrow} \\ \stackrel{S}{\longrightarrow} \\ \stackrel{S}{\longrightarrow} \\ \stackrel{N}{\longrightarrow} \\ \stackrel{N}{\longrightarrow}$$

- and the <substituent group β> represents the following:
  a hydroxyl group, a nitro group, a lower alkanoyl group, a lower alkanoyloxy group, a lower alkoxy group, and a lower alkoxy carbonyl group.
- 4. The compound according to claim 3 or a pharmaceutically acceptable salt thereof, wherein A represents O; the <substituent group  $\alpha>$  represents the following:

the <substituent group  $\beta>$  represents the following:

a hydroxyl group, a lower alkanoyl group, a lower alkanoyloxy group, and a lower alkoxy carbonyl group.

5. The compound according to claim 3 or a pharmaceutically acceptable salt thereof, wherein A represents NH or CH<sub>2</sub>; the <substituent group  $\alpha>$  represents the following:



; and the <substituent group  $\beta$ > represents the following:

- 5 a hydroxyl group, a lower alkanoyl group, and a lower alkyl carbonyloxy group.
  - 6. A compound according to claim 1 which is:
  - 6-N-(2,5-dihydroxymethyl-3-thienylmethyl)amino- $12,13-dihydro-2,10-dihydroxy-12-\beta-D-glucopyranosyl-<math>5H-indolo[2,3-a]$ pyrrolo[3,4-c]carbazole-5,7(6H)-dione;
- 6-N-pyrazinylmethylamino-12,13-dihydro-2,10-dihydroxy-12-β-D-glucopyranosyl-5H-indolo-[2,3-a]pyrrolo[3,4-c]carbazole-5,7(6H)-dione;
  - 6-N-(4-quinolinylmethylamino)-12,13-dihydro-2,10-dihydroxy-12-β-D-glucopyranosyl-5H-indolo[2,3-a]pyrrolo[3,4-c]carbazole-5,7(6H)-dione;
- glucopyranosyl-5H-indolo[2,3-a]pyrrolo[3,4-c]carbazole-5,7(6H)-dione;
  - 6-N-(2-1H-pyrrolylmethyl)amino-12,13-dihydro-2,10-dihydroxy-12-β-D-glucopyranosyl-5H-indolo[2,3-a]pyrrolo[3,4-c]carbazole-5,7(6H)-dione;
  - 6-N-(3-1H-pyrazolylmethyl)amino- $12,13-dihydro-2,10-dihydroxy-<math>12-\beta-D-glucopyranosyl-5H-indolo[2,3-a]$ pyrrolo[3,4-c]carbazole-5,7(6H)-dione;
- 6-N-(4-1H-imidazolylmethyl)amino-12,13-dihydro-2,10-dihydroxy-12-β-D-glucopyranosyl-5H-indolo[2,3-a]pyrrolo[3,4-c]carbazole-5,7(6H)-dione;
  - 6-N-(2-methoxycarbonyl-6-pyridylmethyl)amino-12,13-dihydro-2,10-dihydroxy-12- $\beta$ -D-glucopyranosyl-5H-indolo[2,3-a]pyrrolo[3,4-c]carbazole-5,7(6H)-dione;
  - $6-N-(4-pyridylmethoxy)-12,13-dihydro-2,10-dihydroxy-12-\beta-D-glucopyranosyl-5H-indolo-10-dihydroxy-12-\beta-D-glucopyranosyl-5H-indolo-10-dihydroxy-12-\beta-D-glucopyranosyl-5H-indolo-10-dihydroxy-12-\beta-D-glucopyranosyl-5H-indolo-10-dihydroxy-12-\beta-D-glucopyranosyl-5H-indolo-10-dihydroxy-12-\beta-D-glucopyranosyl-5H-indolo-10-dihydroxy-12-\beta-D-glucopyranosyl-5H-indolo-10-dihydroxy-12-\beta-D-glucopyranosyl-5H-indolo-10-dihydroxy-12-\beta-D-glucopyranosyl-5H-indolo-10-dihydroxy-1$
- 25 [2,3-a]pyrrolo[3,4-c]carbazole-5,7(6H)-dione;
  - 6-N-(6-hydroxymethyl-2-pyridylmethoxy)-12,13-dihydro-2,10-dihydroxy-12-β-D-glucopyranosyl-5H-indolo[2,3-a]pyrrolo[3,4-c]carbazole-5,7(6H)-dione;
  - 6-N-(3-hydroxymethyl-4-pyridylmethoxy)-12,13-dihydro-2,10-dihydroxy-12-β-D-glucopyranosyl-5H-indolo[2,3-a]pyrrolo[3,4-c]carbazole-5,7(6H)-dione; or
- 6-N-(2-(4-pyridyl)ethyl)-12,13-dihydro-2,10-dihydroxy-12-β-D-glucopyranosyl-5H-indolo-[2,3-a]pyrrolo[3,4-c]carbazole-5,7(6H)-dione, or a pharmaceutically acceptable salt thereof.
- 7. A pharmaceutical composition comprising the compound according to any one of claims 1 to 6 as an active ingredient, together with a pharmaceutically acceptable

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carrier or diluent.

8. An antitumor agent comprising the compound according to any one of claims 1 to 6 as an active ingredient, together with a pharmaceutically acceptable carrier or diluent.